

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111422\
 Data File : VX032799.D
 Acq On : 14 Nov 2022 16:59
 Operator : JC/MD
 Sample : N5572-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 31915

Quant Time: Nov 15 05:09:14 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 08:28:06 2022
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.556	168	95736	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	157975	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	143755	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	71290	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	67273	50.573	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.140%
35) Dibromofluoromethane	5.385	113	54429	48.302	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.600%
50) Toluene-d8	8.647	98	195528	49.901	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.800%
62) 4-Bromofluorobenzene	11.079	95	76799	50.176	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.360%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.989	59	5335	27.555	ug/l #	89
16) Acetone	2.385	43	108665	224.758	ug/l	98
18) Methyl Acetate	2.702	43	9616	6.116	ug/l	99
20) Methylene Chloride	2.788	84	851	0.768	ug/l #	92
25) 2-Butanone	4.562	43	162305	216.684	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	14409	9.121	ug/l	98
52) Toluene	8.720	92	1018	0.351	ug/l	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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